Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_E\Data\BE070615\
Data File : BE090056.D
Acq On : 6 Jul 2015 15:01
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
LabSampleId : SSTDCCC001

Quant Time: Jul 07 01:13:18 2015
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\8270-SIM-BE062915.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jun 30 07:01:01 2015
Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
Max. RRF Dev : 25% Max. Rel. Area : 150%

		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	61	-0.01
2		1,4-Dioxane	0.686	0.639	6.9	59	-0.02
3		n-Nitrosodimethylamine	0.930	0.732	21.3	50	-0.02
4	S	2-Fluorophenol	1.149	1.013	11.8	54	-0.01
5	S	Phenol-d6	1.608	1.419	11.8	55	0.00
6	I	Naphthalene-d8	1.000	1.000	0.0	58	-0.01
7	S	Nitrobenzene-d5	0.397	0.339	14.6	50	-0.01
8		Nitrobenzene	0.421	0.344	18.3	49#	0.00
9		Naphthalene	1.130	1.027	9.1	54	-0.01
10		Hexachlorobutadiene	0.178	0.169	5.1	56	-0.01
11		2-Methylnaphthalene	0.755	0.630	16.6	49#	0.00
12	I	Acenaphthene-d10	1.000	1.000	0.0	49#	0.00
13	S	2,4,6-Tribromophenol	0.147	0.093	36.7#	32#	0.00
14	S	2-Fluorobiphenyl	1.503	1.493	0.7	50#	0.00
15		Acenaphthylene	9.019	8.138	9.8	45#	0.00
16		Acenaphthene	1.300	1.152	11.4	44#	-0.01
17		Fluorene	1.768	1.536	13.1	43#	-0.02
18	I	Phenanthrene-d10	1.000	1.000	0.0	42#	-0.02
19		4-Bromophenyl-phenylether	0.213	0.192	9.9	39#	-0.02
20		Hexachlorobenzene	0.885	0.814	8.0	40#	-0.02
21		Pentachlorophenol	0.086	0.315	-266.3#	162#	0.00
22		Phenanthrene	1.281	1.094	14.6	37#	-0.02
23		Anthracene	1.172	0.999	14.8	36#	0.00
24		Fluoranthene	1.431	1.168	18.4	35#	-0.01
25	I	Chrysene-d12	1.000	1.000	0.0	37#	-0.02
26		Pyrene	1.192	1.111	6.8	35#	-0.01
27	S	Terphenyl-d14	0.833	0.771	7.4	34#	-0.01
28		Benzo(a)anthracene	1.249	1.079	13.6	33#	-0.01
29		Chrysene	1.297	1.137	12.3	33#	-0.01
30		Bis(2-ethylhexyl)phthalate	0.398	0.192	51.8#	19#	-0.02
31		Indeno(1,2,3-cd)pyrene	1.243	1.125	9.5	34#	-0.04
32	I	Perylene-d12	1.000	1.000	0.0	35#	-0.02
33		Benzo(b)fluoranthene	1.103	1.105	-0.2	35#	-0.02
34		Benzo(k)fluoranthene	1.232	1.200	2.6	34#	-0.02
35	C	Benzo(a)pyrene	1.019	0.990	2.8	34#	-0.02
36		Dibenzo(a,h)anthracene	1.046	0.979	6.4	33#	-0.04
37		Benzo(g,h,i)perylene	1.060	0.997	5.9	33#	-0.04

(#) = Out of Range SPCC's out = 0 CCC's out = 0